of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 13:53:24 ON 15 JUL 2004

=> le reg
THIS COMMAND NOT AVAILABLE IN THE CURRENT FILE
Some commands only work in certain files. For example, the EXPAND
command can only be used to look at the index in a file which has an
index. Enter "HELP COMMANDS" at an arrow prompt (=>) for a list of
commands which can be used in this file.

=> file reg
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.21 0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 13:53:34 ON 15 JUL 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 13 JUL 2004 HIGHEST RN 709042-93-3 DICTIONARY FILE UPDATES: 13 JUL 2004 HIGHEST RN 709042-93-3

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

Please note that search-term pricing does apply when conducting ${\tt SmartSELECT}$ searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at: http://www.cas.org/ONLINE/DBSS/registryss.html

=>

Uploading C:\Examination Auxillary files\09995987\09995987 third try fixed H.str

chain nodes :

1 2 3 4 5 6 7 8 9 10 11 12 14

chain bonds :

1-2 1-12 2-3 2-7 3-4 3-8 3-9 4-5 5-6 5-10 6-11 6-14

exact/norm bonds:

4-5 5-6 6-11 6-14

exact bonds :

1-12 2-3 3-4 3-8 3-9 5-10

normalized bonds :

1-2 2-7

G1:C,O,S,N

Hydrogen count : 4:>= minimum 2 Match level :

1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS 12:CLASS 14:CLASS

L1STRUCTURE UPLOADED

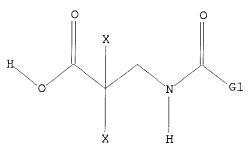
=> id 11

COMBINATION OF STRUCTURE AND TEXT TERMS NOT VALID The query entered contains both search terms created by structure-building or screen commands and text search terms. L#s created via the STRUCTURE or SCREEN commands must be searched in the structures files separately from text terms or profiles. The L# answer sets from structure searches can be used in crossover searches and can be combined with text terms.

=> d 11

L1 HAS NO ANSWERS

L1



G1 C, O, S, N

Structure attributes must be viewed using STN Express query preparation.

=> search l1 sss sam SAMPLE SEARCH INITIATED 13:54:06 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 21 TO ITERATE

100.0% PROCESSED 21 ITERATIONS SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE** 146 TO

PROJECTED ITERATIONS: 694 PROJECTED ANSWERS: OTO 0

0 SEA SSS SAM L1

=> search 11 sss full FULL SEARCH INITIATED 13:54:14 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 508 TO ITERATE

100.0% PROCESSED 508 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

L3 1 SEA SSS FUL L1

=> d scan

L3 1 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

N Propanoic acid, 2,2-difluoro-3-[[(phenylmethoxy)carbonyl]amino]- (9CI)

MF Cl1 H11 F2 N O4

 $\begin{array}{c} \text{O} \\ \parallel \\ \text{Ph-CH}_2 - \text{O-C-NH-CH}_2 - \text{CF}_2 - \text{CO}_2 \text{H} \end{array}$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus
COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 155.84 156.05

FULL ESTIMATED COST

FILE 'CAPLUS' ENTERED AT 13:54:28 ON 15 JUL 2004 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

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FILE COVERS 1907 - 15 Jul 2004 VOL 141 ISS 3 FILE LAST UPDATED: 14 Jul 2004 (20040714/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> 13

L4 1 L3

=> d 14 ti fbib abs

- L4 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN
- TI Preparation of ureidomethylbenzoylaminodifluoropropionates as glucagon antagonists/inverse agonists.
- AN 2002:391685 CAPLUS
- DN 136:385945
- TI Preparation of ureidomethylbenzoylaminodifluoropropionates as glucagon antagonists/inverse agonists.

```
Jorgensen, Anker Steen; Madsen, Peter
PA
     Novo Nordisk A/S, Den.
SO
     PCT Int. Appl., 85 pp.
     CODEN: PIXXD2
DT
     Patent
     English
FAN.CNT 1
     PATENT NO.
                      KIND DATE
                                           APPLICATION NO. DATE
     _____ ____
                                            _____
                                                             _____
PΙ
     WO 2002040446
                      A1
                            20020523
                                           WO 2001-DK760 20011115
         W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN,
             CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR,
             LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PH, PL,
             PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG,
             UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM
         RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH,
             CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR,
             BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG
                                            DK 2000-1733 A 20001117
     AU 2002023502
                       Α5
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                                            AU 2002-23502
                                                              20011115
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     EP 1345891
                                            EP 2001-996529 20011115
                       A1
                             20030924
         R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,
             IE, SI, LT, LV, FI, RO, MK, CY, AL, TR
                                            DK 2000-1733
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                                            WO 2001-DK760 W 20011115
     JP 2004513936
                       Т2
                             20040513
                                            JP 2002-542774
                                                             20011115
                                            DK 2000-1733 A 20001117
                                            WO 2001-DK760 W 20011115
     US 2003027849
                       A1
                             20030206
                                            US 2001-995987
                                                              20011116
                                            DK 2000-1733 A 20001117
                                            US 2000-252322PP 20001120
OS
    MARPAT 136:385945
AΒ
     HO2CCF2CH2NHCOZCHR2N(E)XD [R2 = H, alkyl; Z = (substituted) arylene,
     heteroarylene; X = (CH2)q(CR12R13)r(CH2)s, CO(CR12R13)r(CH2)s,
     NR11CO(CR12R13)r(CH2)s, etc.; r = 0, 1; s = 0-3; R11, R12, R13 = H,
    alkyl; D = (substituted) Ph, naphthyl, pyridyl, indenyl, benzothienyl,
     thienyl, furyl, benzofuryl, etc.; E = (substituted) cyclohexyl, Ph, PhCH2, PhCH2CH2, indanyl, benzhydryl, etc.], were prepared Thus, Me
     4-[(4-cyclohex-1-enylphenylamino)methy1]benzoate (preparation given) in CH2Cl2
     containing diisopropylethylamine was treated with 3,5-dichlorophenyl
     isocyanate to give a residue which was saponified with LiOH. The resulting
     acid in DMF was treated with 3-[(dimethyliminium)(dimethylamino)methyl]-
     1,2,3-benzotriazol-1-ium-1-olate hexafluorophosphate,
     diisopropylethylamine, Me 3-amino-2,2-difluoropropionate hydrochloride to
     give the uncharacterized amide ester, which was saponified with aqueous LiOH in
     THF/MeOH to give 3-[4-[1-(4-cyclohex-1-enylphenyl)-3-(3,5-
     dichlorophenyl)ureidomethyl]benzoylamino]-2,2-difluoropropionic acid. In
     a human glucagon receptor binding assay, title compds. showed IC50<1000
    nM.
RE.CNT 3
              THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD
              ALL CITATIONS AVAILABLE IN THE RE FORMAT
=> logoff hold
COST IN U.S. DOLLARS
                                                  SINCE FILE
                                                                   TOTAL
                                                        ENTRY
                                                                 SESSION
FULL ESTIMATED COST
                                                        3.14
                                                                  159.19
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)
```

SINCE FILE

ENTRY

TOTAL

SESSION

ΙN

-0.74 -0.74CA SUBSCRIBER PRICE

SESSION WILL BE HELD FOR 60 MINUTES STN INTERNATIONAL SESSION SUSPENDED AT 13:55:10 ON 15 JUL 2004

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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NEWS 2 "Ask CAS" for self-help around the clock

NEWS 3 May 12 EXTEND option available in structure searching
NEWS 4 May 12 Polymer links for the POLYLINK command completed in REGISTRY
NEWS 5 May 27 New UPM (Update Code Maximum) field for more efficient patent SDIs in CAplus

NEWS 6 May 27 CAplus super roles and document types searchable in REGISTRY

NEWS $\,$ 7 Jun 22 STN Patent Forums to be held July 19-22, 2004

NEWS 8 Jun 28 Additional enzyme-catalyzed reactions added to CASREACT

NEWS 9 Jun 28 ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG, and WATER from CSA now available on STN(R)

NEWS 10 Jul 12 BEILSTEIN enhanced with new display and select options, resulting in a closer connection to BABS

NEWS EXPRESS MARCH 31 CURRENT WINDOWS VERSION IS V7.00A, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 26 APRIL 2004

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NEWS WWW CAS World Wide Web Site (general information)

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FILE 'HOME' ENTERED AT 06:14:34 ON 16 JUL 2004

=> ile req

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